

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN032619\
 Data File : BN004973.D
 Acq On : 26 Mar 2019 13:06
 Operator : JU/SJ
 Sample : SSTDICC020
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDICC020

Quant Time: Mar 26 15:15:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN032619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 15:08:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	9334	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	42182	20.00	ng	0.00
39) Acenaphthene-d10	14.32	164	25011	20.00	ng	0.00
64) Phenanthrene-d10	17.07	188	56062	20.00	ng	0.00
76) Chrysene-d12	21.27	240	56920	20.00	ng	0.00
87) Perylene-d12	23.52	264	51808	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.27	112	21693	41.92	ng	0.00
7) Phenol-d6	6.85	99	29071	40.50	ng	0.00
23) Nitrobenzene-d5	8.84	82	28108	41.94	ng	0.00
42) 2,4,6-Tribromophenol	15.82	330	16251	43.59	ng	0.00
45) 2-Fluorobiphenyl	12.94	172	76298	40.72	ng	0.00
79) Terphenyl-d14	19.70	244	122101	42.23	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.21	88	3906	19.527	ng	97
3) Pyridine	3.60	79	11112	18.582	ng	99
4) n-Nitrosodimethylamine	3.53	42	3292	18.043	ng	# 94
6) Aniline	7.02	93	18463	20.815	ng	99
8) 2-Chlorophenol	7.25	128	13865	21.166	ng	98
9) Benzaldehyde	6.83	77	9329	18.991	ng	97
10) Phenol	6.88	94	15417	19.908	ng	99
11) bis(2-Chloroethyl)ether	7.11	93	11912	19.401	ng	97
12) 1,3-Dichlorobenzene	7.56	146	14971	20.257	ng	98
13) 1,4-Dichlorobenzene	7.71	146	15331	20.297	ng	99
14) 1,2-Dichlorobenzene	8.02	146	14678	20.091	ng	99
15) Benzyl Alcohol	7.92	79	10786	20.163	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.20	45	14668	18.629	ng	99
17) 2-Methylphenol	8.12	107	10945	20.482	ng	99
18) Hexachloroethane	8.74	117	5380	21.885	ng	97
19) n-Nitroso-di-n-propylamine	8.48	70	9725	20.125	ng	98
20) 3+4-Methylphenols	8.45	107	14829	19.928	ng	99
22) Acetophenone	8.50	105	19630	20.411	ng	100
24) Nitrobenzene	8.88	77	14184	20.626	ng	99
25) Isophorone	9.39	82	27842	23.992	ng	99
26) 2-Nitrophenol	9.59	139	8081	20.669	ng	98
27) 2,4-Dimethylphenol	9.65	122	11563	20.692	ng	99
28) bis(2-Chloroethoxy)methane	9.88	93	17186	21.271	ng	98
29) 2,4-Dichlorophenol	10.12	162	13110	19.938	ng	99
30) 1,2,4-Trichlorobenzene	10.32	180	14458	19.292	ng	99
31) Naphthalene	10.51	128	44819	21.828	ng	100
32) Benzoic acid	9.78	122	7612	18.425	ng	98
33) 4-Chloroaniline	10.63	127	18272	19.944	ng	100
34) Hexachlorobutadiene	10.78	225	9308	20.742	ng	100
35) Caprolactam	11.41	113	4339	18.714	ng	98
36) 4-Chloro-3-methylphenol	11.76	107	14397	20.669	ng	99
37) 2-Methylnaphthalene	12.13	142	31974	21.804	ng	100
38) 1-Methylnaphthalene	12.35	142	31490	22.170	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.50	216	17380	19.331	ng	99

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41) Hexachlorocyclopentadiene	12.46	237	8012	20.672	nq	98
43) 2,4,6-Trichlorophenol	12.75	196	10825	19.391	nq	99
44) 2,4,5-Trichlorophenol	12.83	196	11841	19.885	nq	100
46) 1,1'-Biphenyl	13.15	154	42303	19.097	nq	100
47) 2-Chloronaphthalene	13.19	162	32414	19.448	nq	99
48) 2-Nitroaniline	13.42	65	8617	20.750	nq	99
49) Acenaphthylene	14.05	152	55097	22.408	nq	100
50) Dimethylphthalate	13.79	163	41873	21.371	nq	99
51) 2,6-Dinitrotoluene	13.92	165	9476	20.719	nq	96
52) Acenaphthene	14.39	154	31005	20.792	nq	99
53) 3-Nitroaniline	14.25	138	9547	19.526	nq	95
54) 2,4-Dinitrophenol	14.46	184	4326	23.351	nq	96
55) Dibenzofuran	14.73	168	49892	20.311	nq	100
56) 4-Nitrophenol	14.57	139	6962	18.636	nq	98
57) 2,4-Dinitrotoluene	14.70	165	12938	20.929	nq	# 98
58) Fluorene	15.37	166	40666	22.404	nq	100
59) 2,3,4,6-Tetrachlorophenol	14.96	232	11066	20.590	nq	# 99
60) Diethylphthalate	15.15	149	42605	20.551	nq	99
61) 4-Chlorophenyl-phenylether	15.37	204	21021	19.672	nq	99
62) 4-Nitroaniline	15.42	138	9746	19.863	nq	99
63) Azobenzene	15.66	77	35127	20.759	nq	100
65) 4,6-Dinitro-2-methylphenol	15.47	198	6928	21.077	nq	97
66) n-Nitrosodiphenylamine	15.59	169	35623	20.061	nq	99
67) 4-Bromophenyl-phenylether	16.26	248	13823	19.548	nq	97
68) Hexachlorobenzene	16.37	284	16074	19.531	nq	100
69) Atrazine	16.54	200	13072	19.680	nq	100
70) Pentachlorophenol	16.73	266	7312	13.330	nq	98
71) Phenanthrene	17.12	178	63831	21.795	nq	100
72) Anthracene	17.20	178	63394	22.145	nq	99
73) Carbazole	17.49	167	58788	20.925	nq	100
74) Di-n-butylphthalate	18.03	149	72206	21.781	nq	100
75) Fluoranthene	19.14	202	74874	23.716	nq	99
77) Benzidine	19.33	184	39036	19.831	nq	99
78) Pyrene	19.50	202	76373	19.878	nq	99
80) Butylbenzylphthalate	20.40	149	31423	19.083	nq	100
81) Benzo(a)anthracene	21.26	228	71583	22.108	nq	99
82) 3,3'-Dichlorobenzidine	21.19	252	27762	20.573	nq	98
83) Chrysene	21.31	228	68882	22.244	nq	98
84) Bis(2-ethylhexyl)phthalate	21.17	149	43338	18.149	nq	99
85) Di-n-octyl phthalate	22.05	149	68385	19.282	nq	99
86) Indeno(1,2,3-cd)pyrene	25.77	276	67425	20.625	nq	100
88) Benzo(b)fluoranthene	22.84	252	65981	22.875	nq	100
89) Benzo(k)fluoranthene	22.89	252	60229	20.984	nq	99
90) Benzo(a)pyrene	23.42	252	59097	22.113	nq	99
91) Dibenzo(a,h)anthracene	25.78	278	55479	20.385	nq	99
92) Benzo(g,h,i)perylene	26.47	276	54275	20.508	nq	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

